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REPORT NUMBER 6

ION PHENOMENA

Final Report

by

Robert W. Kiser

May 1968

Physical Research Laboratory
RESEARCH LABORATORIES
Edgewood Arsenal, Maryland 21010

Contract DA-18-035-AMC-718(A)

KANSAS STATE UNIVERSITY
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FOREWORD

The work described in this report was authorized under Project 1C622401A102, Chemical Agent Warning and Detection Techniques (U). This work was started 9 December 1965 and completed 31 July 1967. The experimental data are contained in notebooks KSU-182311-DLD-6 to KSU-182311-DLD-11 and KSU-182311-RES-3 to KSU-182311-RES-5.

Reference to a company and/or product name in this report is only for purposes of information and does not imply approval or recommendation of the product to the exclusion of others, which may also be suitable.

The information in this report has not been cleared for release to the general public.

Acknowledgments

Donald L. Dugger and Richard E. Sullivan have been instrumental in obtaining the experimental findings reported herein.

DIGEST

Mass spectrometric studies of trimethyl phosphorothionate, triphenyl phosphite, diallyl phosphite, diallyl allylphosphonate, triallyl phosphite, triallyl phosphate, and diphenyl phosphite were made to obtain more fundamental information for use in determining the ionization and dissociative ionization processes that occur upon electron impact with these organophosphorus compounds. Mass spectra, appearance potentials, clastograms and metastable transitions obtained in these studies are reported. These results are discussed in relation to a better understanding of the fragmentation pathways in this class of compounds. A summary table of all results determined under this contract is included, and a brief catalog, with literature references, of the mass spectra of organophosphorus compounds summarizes much of the available literature. Future correlative activities are suggested.

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ION PHENOMENA

1. STATEMENT OF PROBLEM

This work was initiated in order (1) to obtain mass spectra, (2) to make energetic studies by appearance potentials, and clastogram determinations, (3) to utilize metastable transition measurements in determining fragmentation pathways, and (4) to form negative parent molecule ions by bombardment of biologically active compounds containing P, N, and S by means of a polonium-210 alpha particle source.

2. BACKGROUND

Preliminary investigations of organophosphorus compounds such as GB and GD have indicated that these biologically active materials are ionized at low potentials (ca. 11.5 eV) and that at electron energies above approximately 15 eV there is significant fragmentation of the parent positive ion. However, the ionization and dissociative ionization processes are poorly understood, attributable largely to a lack of fundamental information concerning similar processes in simpler organophosphorus compounds. Because little or no information is available about the processes in the biologically active compounds, it is necessary to study both the simpler and the more complex organophosphorus molecules.

The study of both the positive and negative ions formed in the ionization and dissociation processes is important to understanding the fragmentation processes that occur. It has been reported that the G agents capture low energy electrons to form parent negative ions, but almost certainly other negatively charged fragment ions will be formed with higher energy electrons through dissociative electron attachment. No information is available about negative ion clastograms that would indicate possible routes of fragmentation of these species.

3. APPROACH TO THE PROBLEM

3.1 Materials

The materials to be studied included trimethyl phosphorothionate, triphenyl phosphite, diallyl phosphite, diallyl allylphosphonate, triallyl phosphite, triallyl phosphate, and diphenyl phosphite

3.2. Equipment

The Bendix time-of-flight mass spectrometer, together with its various modifications, has been described in previous reports.[1-8]

3.3. Techniques of Study

The techniques of study have been outlined in detail in earlier reports.[4-8]

4. RESULTS

4.1. Trimethyl Phosphorothionate

Appearance potentials for twelve ions formed by electron impact from trimethyl phosphorothionate are given in Table 1. Processes for the formation of five ions are suggested, based on the energetics and clastogram data.

4.2. Triphenyl Phosphite

The 70 eV mass spectrum of triphenyl phosphite was determined and is shown in Figure 1. The clastogram for this molecule is shown in Figures 2 and 3. The appearance potentials listed in Table 2 indicate that the major ions in the mass spectrum of this compound are due to phenol, probably formed by thermal decomposition of the sample. Thus, the base peak at m/q = 94 has an appearance potential of 8.8 ±0.2 eV as determined by electron impact, and the photoionization value for the ionization potential of phenol is known to be 8.5 eV.[9] This comparison and the other data of Table 2 strongly suggest that the majority of the ions below m/q = 94 are due to phenol. The small quantities of ions at m/q = 149, 156 and 167 precluded appearance potential measurements.

4.3. Diallyl Phosphite

The 70 ev mass spectrum of diallyl phosphite was given in Quarterly Progress Report No. 5.[8] The clastogram and several preliminary appearance potential measurements for ions formed from this compound by electron impact were given also in that report.[8] Table 3 given in the Appendix presents a more complete treatment of the previous data.

4.4. Diallyl Allylphosphonate

In Table 4 are listed appearance potentials for thirteen ions formed from diallyl allylphosphonate. However, only for four ions have probable processes of formation been indicated. The ions of m/q = 79, 80 and 81 may be due to PO_3^+, HPO_3^+ and $H_2PO_3^+$, but the appearance potentials do not support this. It will be necessary in

future studies to utilize high resolution mass spectrometry in order to positively identify these ions and thereby permit the process assignments to be made from the energetics and observed metastable transitions.

4.5. Triallyl Phosphite

Even as with diallyl allylphosphonate, the m/q = 79, 80 and 81 ions are not established with sufficient certainty to permit the assignment of probable processes for these ions. Four ions that involve simpler fragmentation have been assigned probable processes in Table 5. Appearance potentials for twelve ions have been determined.

4.6. Triallyl Phosphate

Table 6 lists the principal ions in the 70 eV positive ion mass spectrum together with appearance potentials determined for six different ions using the energy compensation method. The energetic data have been employed in assigning probable processes wherever possible. However, a lack of certain knowledge of composition of many ions does not permit a more detailed analysis of the fragmentation processes at this time. The logarithmic clastogram for this compound is given in Figures 4 and 5.

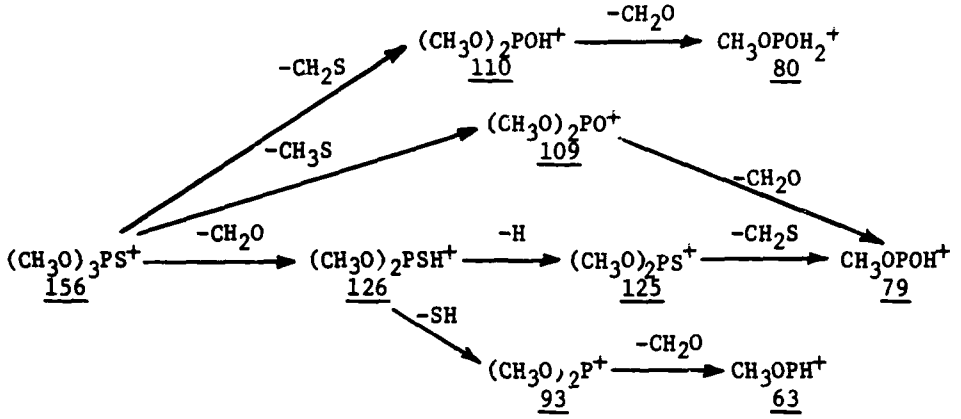
4.7. Diphenyl Phosphite

The 70 eV mass spectrum, the clastogram, and some preliminary appearance potentials for diphenyl phosphite were given previously in Quarterly Progress Report No. 5 [8]. Table 7 presents evidence that the majority of the mass spectrum derives from phenol, even as was noted for triphenyl phosphite. Apparently this is due to thermal decomposition of these materials; this should be studied further in future efforts.

5. DISCUSSION

5.1. Trimethyl Phosphorothionate

Utilizing the metastable transition data, the results from the appearance potential determinations, and the clastograms obtained, the following partial fragmentation scheme may be constructed.



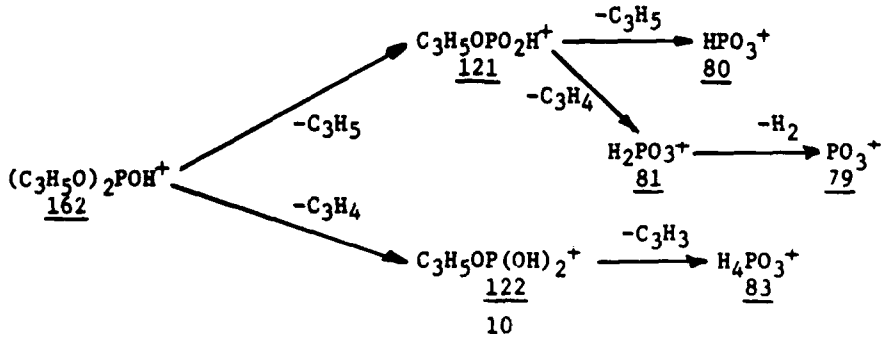
From the metastable transition data, the second process in Table 1 for the formation of the m/q = 79 ion is preferred; thus, the value of $\Delta H_f(79^+)$ = 126 kcal/mole is also preferred. This indicates that the m/q = 79 ion is predominantly CH_3OPOH^+. The unique fragmentations involving the retention and loss of the sulfur atom suggest either that the CH_3 group is very mobile or that rapid equilibria may be established between isomeric forms of these gaseous ionic species.

5.2. Triphenyl Phosphite

In that the mass spectrum and appearance potentials determined suggest that phenol may be formed from this compound through thermal decomposition, and that the present observations do not differentiate any phenol from the ions formed directly from the title compound by electron bombardment, no further discussion of the present results is merited.

5.3. Diallyl Phosphite

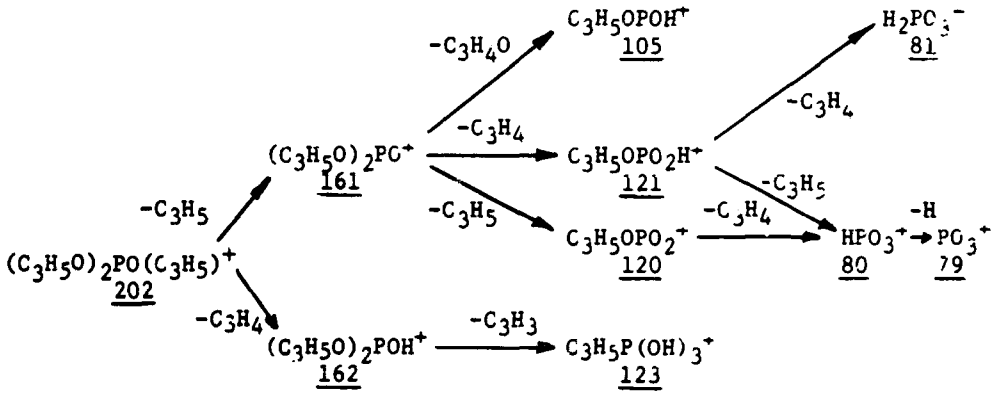
The absence of the parent ion, and the very low appearance potential for the lowest energetic requirement of the fragment ions suggests that the ionization potential of the parent molecule is approximately 10.0 eV. From the clastogram data and the appearance potentials determined in this work, the following partial fragmentation scheme has been constructed:



It should be noted that this scheme for the fragmentation of ions produced in the mass spectrum of diallyl phosphite is significantly different from that originally suggested in Quarterly Progress Report No. 5. (8)

5.4. Diallyl Allylphosphonate

The following fragmentation scheme for the unimolecular decomposition of the diallyl allylphosphonate molecular ion represents a modification of that proposed earlier: (8)



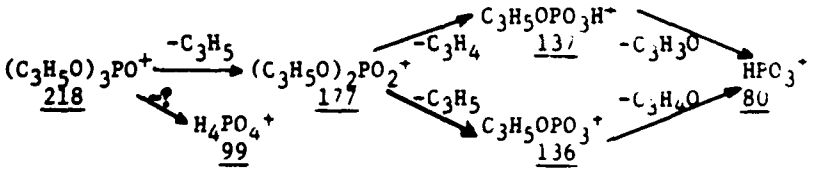
The lowest appearance potential measured suggests that the ionization potential of diallyl allylphosphonate is approximately 10.0 ev.

5.5. Triallyl Phosphite

The additional studies of this molecule have caused no revision in the fragmentation scheme proposed in Quarterly Progress Report No. 5. (8) The ionization potential is estimated to be approximately 8.8 ev, as determined from the lowest appearance potential measured in these studies.

5.6. Triallyl Phosphate

From the acquired appearance potentials and metastable data, the following fragmentation scheme is proposed:



From the appearance potential data, it is suggested that the ionization potential of triallyl phosphate is approximately 9.8 eV.

5.7. Diphenyl Phosphite

Even as with the triphenyl phosphite, the mass spectrum and appearance potentials suggest that phenol may be formed from this compound through thermal decomposition. Therefore, no further discussion of the present results is merited.

6. CONCLUSIONS

It is suggested that future work concentrate on (1) negative ions formed from the various organophosphorus compounds, (2) high resolution mass spectra and precise mass measurements to permit the identification of many of the positive ions formed from these compounds, and (3) detailed metastable transition studies to establish more fully the fragmentation pathways occurring in these compounds. All of these data, together with that obtained under the present contract effort, should yield a reasonably satisfactory body of information capable of providing further insights from correlative studies of the data.

Included in Table 8 in the Appendix is a complete summary of all the findings obtained in these studies. This information, taken together with that summarized in Table 9 and the various other literature-reported results, will provide future investigators with much of the needed material. It is easily concluded that even the data of Tables 8 and 9 are not complete and that additional work will be necessary even in the correlative efforts.

All remaining quantities of the G agents, EA materials, and related substances have been destroyed by hydrolysis in strongly alkaline Chlorox and 15% KOH in 50/50 v/v ethanol-water solutions. HD was similarly treated, except that only Chlorox was employed. All materials and equipment, including the storage refrigerator, have been decontaminated so that no safety hazard remains. Gas masks and atropine sulfate syringes have been stored subject to recall by Edgewood Arsenal.

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APPENDIX

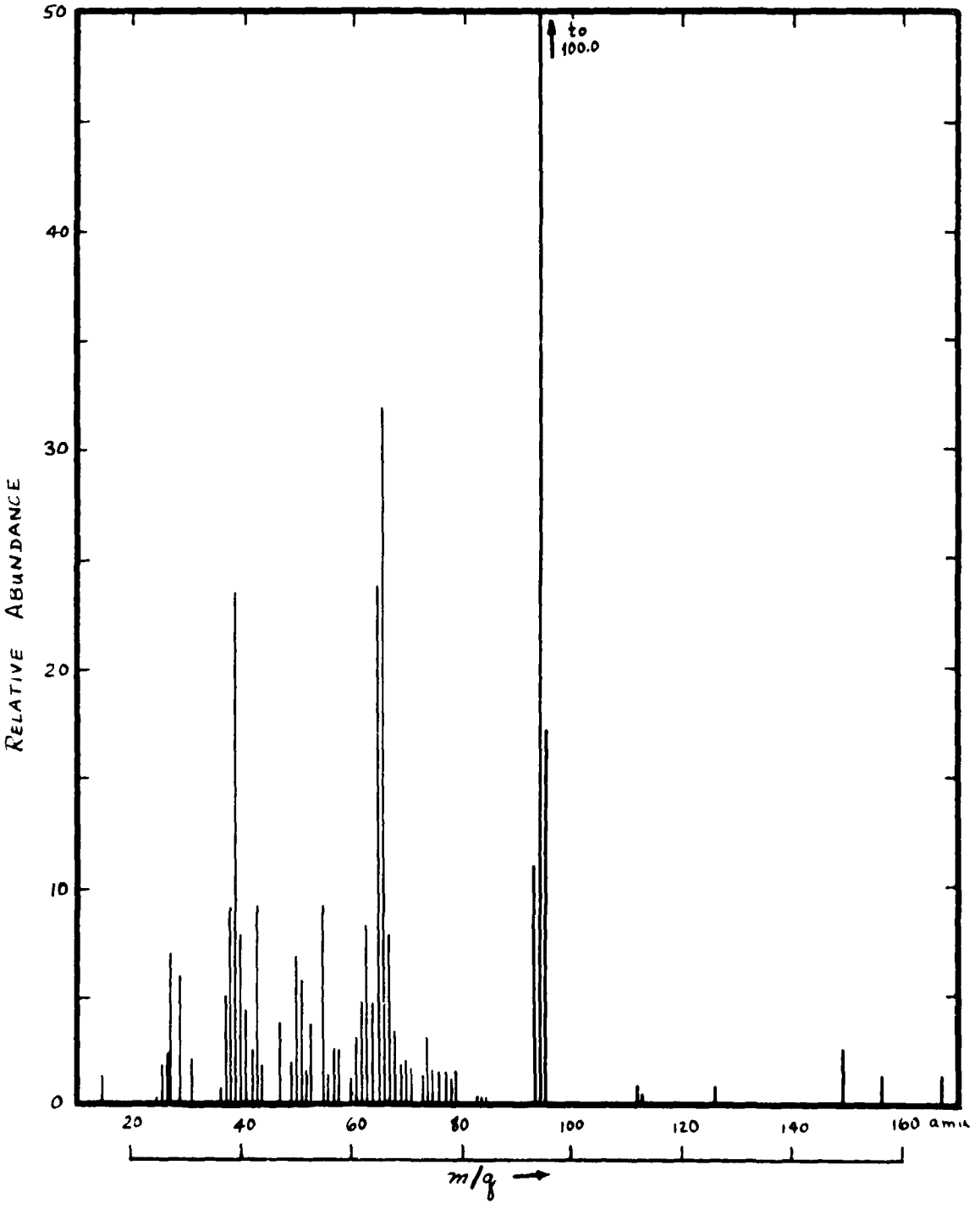
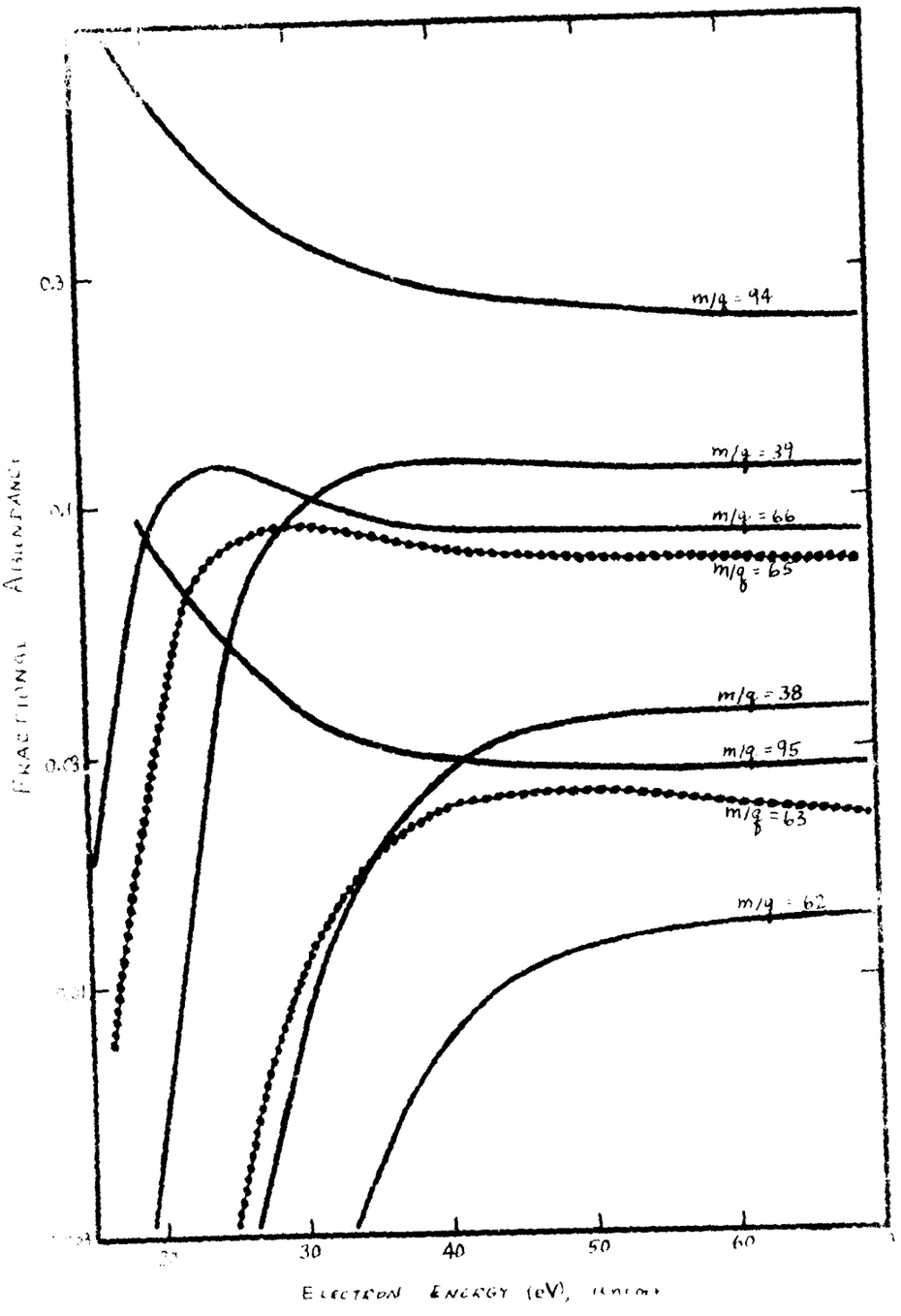


Figure 1. Mass Spectrum of Triphenyl Phosphite, (70 eV).


0.3
0.1
0.03
0.01
m/q = 94
m/q = 39
m/q = 66
m/q = 65
m/q = 38
m/q = 95
m/q = 63
m/q = 62
30
40
50
60
FRACTIONAL ABUNDANCE
ELECTRON ENERGY (eV), (UNCOR

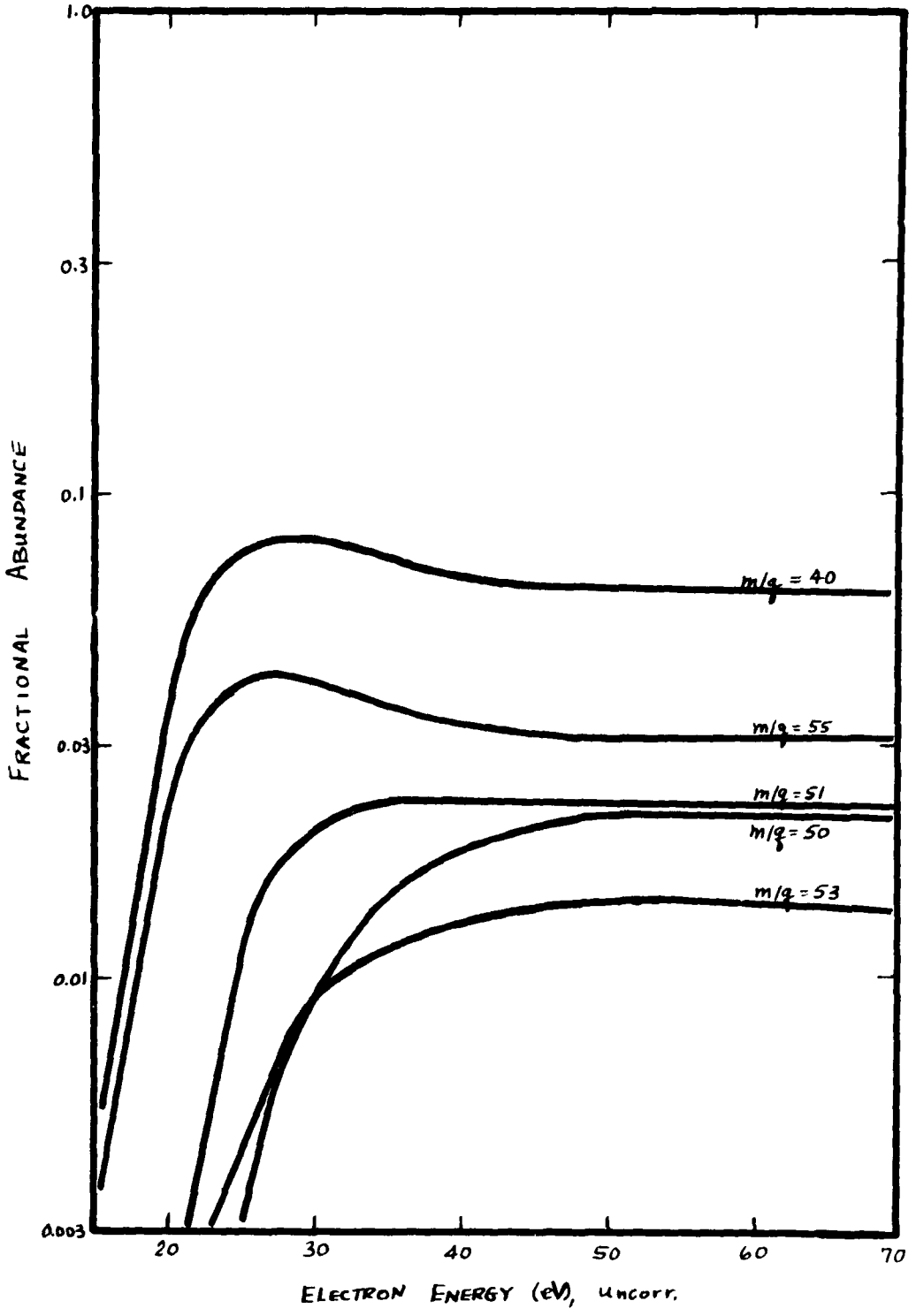


Figure 3. Logarithmic Clastogram for Triphenyl Phosphite [illegible]

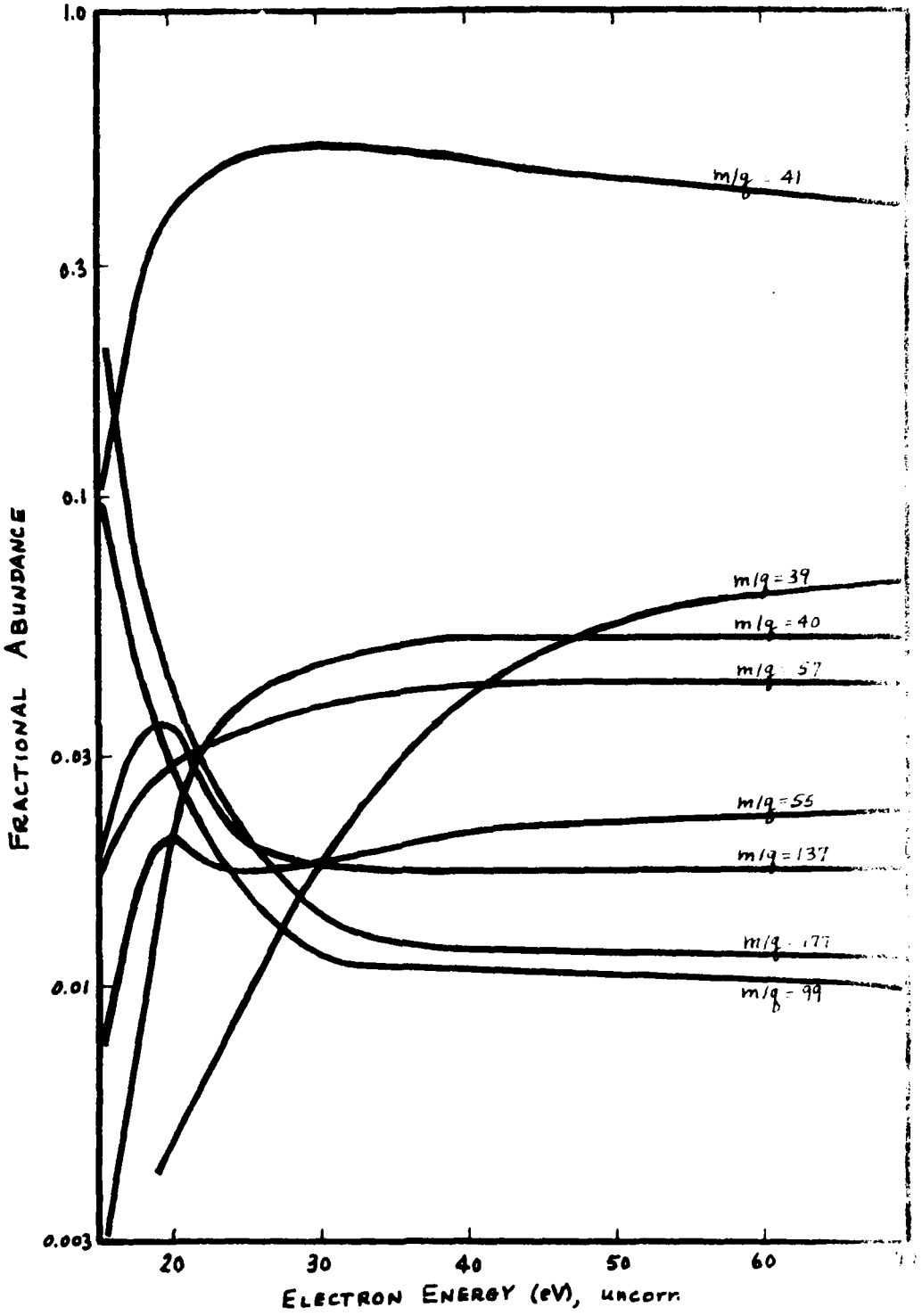


Figure 4. Logarithmic Ebastogram for Triallyl Borate

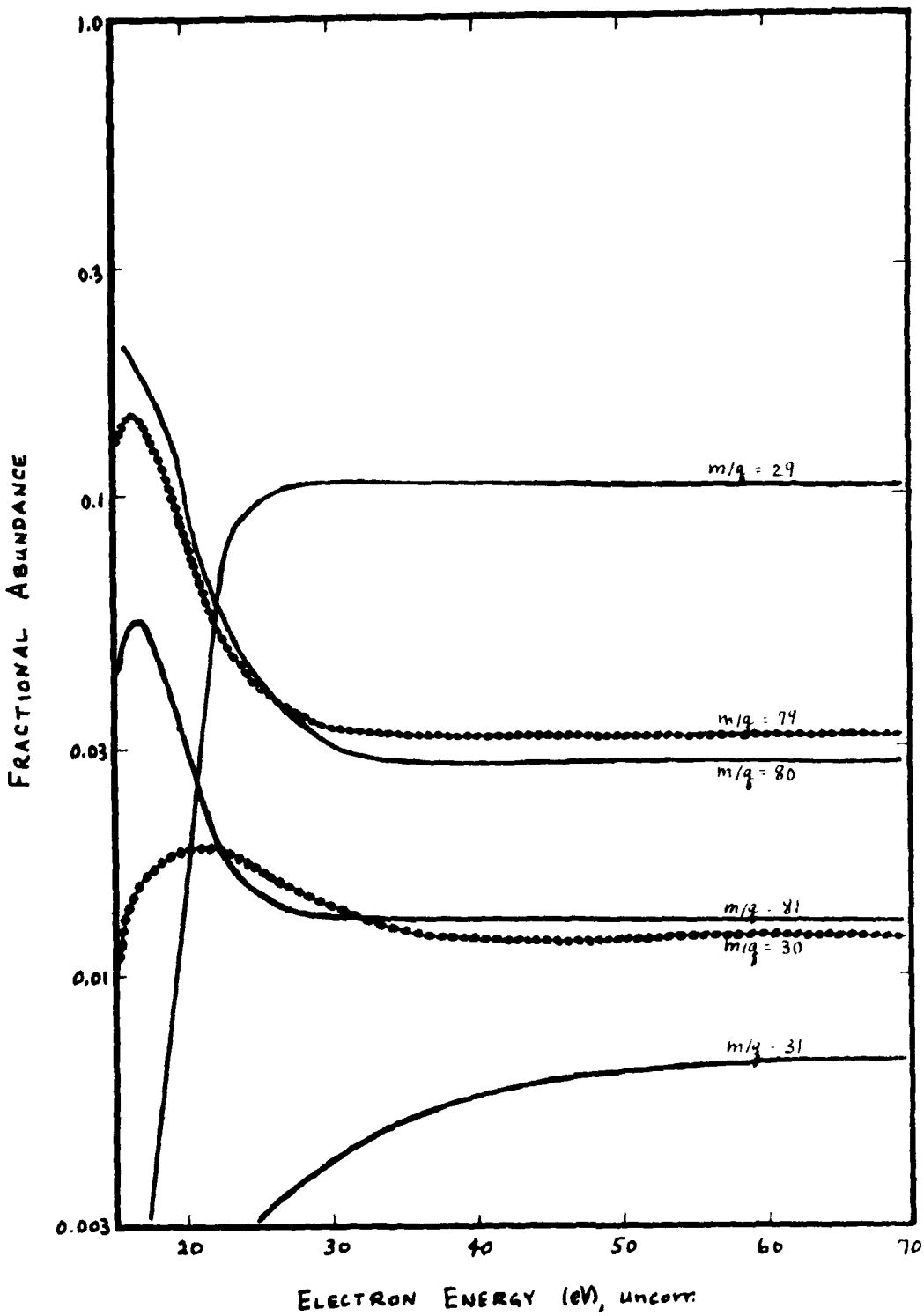

1.0
0.3
0.1
0.03
0.01
0.003
FRACTIONAL ABUNDANCE
m/q = 29
m/q = 74
m/q = 80
m/q = 81
m/q = 30
m/q = 31
20
30
40
50
60
70
ELECTRON ENERGY (eV), uncorr.

Figure 5. [illegible]

Table 1. Mass Spectrum and Appearance Potentials of the Principal Positive Ions formed from Trimethyl Phosphorothionate.[a]

m/e	70 eV Relative Abundance	Appearance Potential (eV)	Probable Process	ΔH_f(ion) (kcal/mole)
156	27.5	[illegible] ±0.2	$SP(OCH_3)_3 \rightarrow SP(OCH_3)_3^+$	8
126	18.5	11.4 ±0.2	$\rightarrow HSP(OCH_3)_2^+ + CH_2O$	97
125	10.0	13.3 ±0.4	$\rightarrow SP(OCH_3)_2^+ + CH_2O + H$	78
93	100.0	12.8 ±0.3	$\rightarrow P(OCH_3)_2^+ + CH_2O + SH$	88
79	11.0	15.1 ±0.3	$\rightarrow PO_3^+ + 2CH_3 + CH_3S$	40
			$\rightarrow CH_3OPOH^+ + CH_2S + CH_3O$	126
63	17.0	16.2 ±0.3	$\rightarrow PS^+ + ?$	
47	19.6	17.3 ±0.3	$\rightarrow PO^+ + ?$	
45	6.4	15.6 ±0.3	$\rightarrow CHS^+ + ?$	
31	14.7	16.0 ±0.3	$\rightarrow OCH_3^+ + ?$	
30	3.2	13.1 ±0.2	$\rightarrow CH_2O^+ + ?$	
29	11.3	17.0 ±0.3	$\rightarrow CHO^+ + ?$	
15	4.8	18.9 ±0.5	$\rightarrow CH_3^+ + ?$	

[illegible] ΔH_f[illegible] kcal/mole.

Table 2. Mass Spectrum and Appearance Potentials for the Principal Positive Ions Observed from Triphenyl Phosphite.[a]

m/q	70 eV Relative Abundance	Appearance Potential (eV)	Probable Process	ΔH_f(ion) (kcal/mole)
167	1.4			
156	1.4			
149	2.6			
94	100.0	8.8 ±0.2[b]	$C_6H_5OH \rightarrow C_6H_5OH^+$	
66	32.0	13.1 ±0.2	$\rightarrow C_5H_6^+ + CO$	
65	23.7	14.4 ±0.2	$\rightarrow C_5H_5^+ + CO + H$	
40	7.8	15.2 ±0.2	$\rightarrow C_3H_4^+ + ?$	
39	23.3	17.4 ±0.2	$\rightarrow C_3H_3^+ + ?$	

a) Estimated $\Delta H_f[P(OC_6H_5)_3] = -78$ kcal/mole.

b) Known ionization potential of phenol (by photoionization) is 8.50 eV.[9]

Table 3. Mass Spectrum and Appearance Potentials of the Principal Positive Ions formed from Diallyl Phosphite.[a]

m/q	70 eV Relative Abundance	Appearance Potential (eV)	Probable Process	ΔH_f(ion) (kcal/mole)
121	36.5	10.6 ±0.2	$(C_3H_5O)_2POH \rightarrow C_3H_5OPO_2H^+ + C_3H_5$	59
83	8.7	11.0 ±0.3	$\rightarrow H_4PO_3^+ + C_3H_3 + C_3H_4$	-28
80	22.8	10.4 ±0.2	$\rightarrow HOPO_2^+ + 2C_3H_5$	23
79	9.5	11.2 ±0.2	$\rightarrow PO_3^+ + C_3H_5 + C_3H_4 + H_2$	28
58	25.0	11.0 ±0.5	$\rightarrow C_3H_5OH^+ + ?$	
57	69.8	11.5 ±0.3	$\rightarrow C_3H_5O^+ + ?$	
55	10.9	11.7 ±0.2	$\rightarrow C_3H_3O^+ + ?$	
41	100.0	13.9 ±0.2	$\rightarrow C_3H_5^+ + ?$	
39	54.5	16.2 ±0.3	$\rightarrow C_3H_3^+ + ?$	

a) Estimated $\Delta H_f\left[(C_3H_5O)_2POH\right]$ = -153 kcal/mole.

Table 4. Mass Spectrum and Appearance Potentials of the Principal Positive Ions formed from Diallyl Allylphosphonate.[a]

m/e	70 eV Relative Abundance	Appearance Potential (eV)		Probable Process	ΔH_f(ion) (kcal/mole)
161	10.8	10.4	±0.2	$(C_3H_5O)_2PO(C_3H_5) \rightarrow C_3H_5OPO_2(C_3H_5)^+ + C_3H_5$	70
123	5.0	10.9	±0.2	$\rightarrow C_3H_5P(OH)_3^+ + C_3H_3 + C_3H_4$	-16
121	13.3	11.0	±0.2	$\rightarrow PO_2(OH)(C_3H_5)^+ + C_3H_4 + C_3H_5$	39
1[illegible]	[illegible].2	12.2	±0.2	$\rightarrow C_3H_5PO_2H^+ + C_3H_5 + C_3H_4O$	134
[illegible]1	[illegible].5	11.6	±0.2	$\rightarrow$?	
[illegible]3	[illegible].5	11.1	±0.2	$\rightarrow$?	
[illegible]	[illegible].2	11.9	±0.2	$\rightarrow$?	
[illegible]	[illegible].3	12.5	±0.2	$\rightarrow H_4PO_2^+ + ?$	
[illegible]	16.5	11.0	±0.2	$\rightarrow C_3H_5OH^+ + ?$	
[illegible]	[illegible].1	12.5	±0.3	$\rightarrow C_3H_5O^+ + ?$	
[illegible]	[illegible].5	13.8	±0.2	$\rightarrow C_3H_4O^+ + ?$	
[illegible]	100.0	14.6	±0.5	$\rightarrow C_3H_5^+ + ?$	
[illegible]	[illegible].2	17.9	±0.3	$\rightarrow C_3H_3^+ + ?$	

[a] Estimated $\Delta H_f\left[(C_3H_5O)_2PO(C_3H_5)\right]$ = -138 kcal/mole.

Table 5. Mass Spectrum and Appearance Potentials of the Principal Positive Ions formed from Triallyl Phosphite.[a]

m/q	70 eV Relative Abundance	Appearance Potential (eV)	Probable Process	ΔH_f(ion) (kcal/mole)
161	4.3	$9.1_5 \pm 0.1_5$	$(C_3H_5O)_3P \rightarrow (C_3H_5O)_2PO^+ + C_3H_5$	65
123	1.8	10.1 ±0.2	$\rightarrow C_3H_5OPO_2H_3^+ + C_3H_3 + C_3H_4$	-10
121	4.0	10.5 ±0.2	$\rightarrow C_3H_5PO_3H^+ + C_3H_5 + C_3H_4$	51
105	3.3	11.0 ±0.2	$\rightarrow C_3H_5OPOH^+ + C_3H_4O + C_3H_5$	131
81	14.2	10.8 ±0.2	$\rightarrow$?	
80	9.5	10.2 ±0.2	$\rightarrow$?	
79	13.1	10.4 ±0.2	$\rightarrow$?	
67	5.9	11.2 ±0.2	$\rightarrow H_4PO_2^+ + ?$	
57	9.0	11.9 ±0.2	$\rightarrow C_3H_5O^+ + ?$	
55	5.1	12.2 ±0.2	$\rightarrow C_3H_3O^+ + ?$	
42	6.3	11.9 ±0.2	$\rightarrow C_2H_2O^+ + ?$	
41	100.0	13.3 ±0.2	$\rightarrow C_3H_5^+ + ?$	
39	21.3		$\rightarrow C_3H_3^+ + ?$	

a) Estimated $\Delta H_f\left[(C_3H_5O)_3P\right] = -114$ kcal/mole.

Table 6. Mass Spectrum and Appearance Potentials for the Principal Positive Ions formed From Triallyl Phosphate.[a]

m/q	70 eV Relative Abundance	Appearance Potential (eV)	Probable Process	ΔH_f(ion) (kcal/mole)
177	2.8	10.2 ±0.3	$(C_3H_5O)_3PO \rightarrow (C_3H_5O)_2PO_2^+ + C_3H_5$	15
137	4.4	12.2 ±0.3	$\rightarrow C_3H_5OPO_3H^+ + C_3H_4 + C_3H_5$	16
93	3.1			
91	3.2			
80	6.2	10.9 ±0.2	$\rightarrow HOPO_2^+ + C_3H_4O + 2C_3H_5$	22
79	7.9			
57	13.3	11.7 ±0.2	$\rightarrow C_3H_5O^+ + ?$	
45	4.6			
41	100.0	11.3 ±0.2	$\rightarrow C_3H_5^+ + ?$	
40	12.7		$\rightarrow C_3H_4^+ + ?$	
39	16.5	15.0 ±0.3	$\rightarrow C_3H_3^+ + ?$	
31	1.7			
29	4.9			
27	16.4			

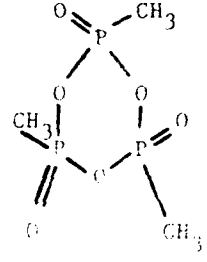
[illegible] ΔH_f[illegible] [illegible] kcal/mole.

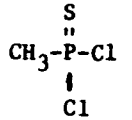
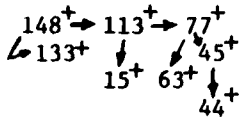
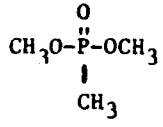
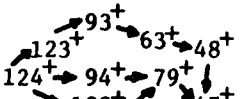
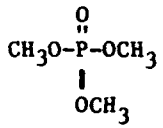
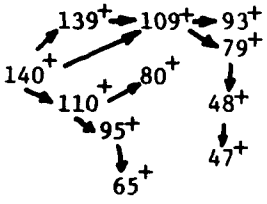
Table 7. Mass Spectrum and Appearance Potentials of the Principal Positive Ions Formed from Diphenyl Phosphite.[a]

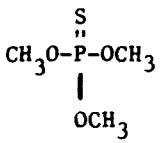
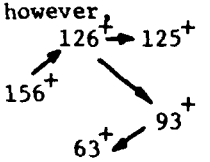
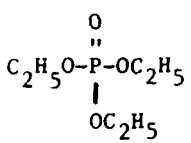
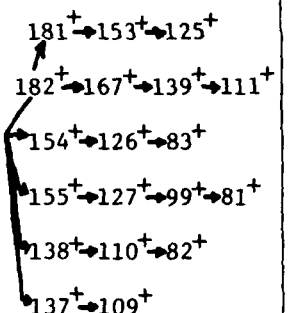
m/q	70 eV Relative Abundance	Appearence Potential (eV)	Probable Process	ΔH_f(ion) (kcal/mole)
167	0.3			
156	0.6			
149	1.4			
95	19.4			
94	100.0	8.5 ±0.2	$C_6H_5OH \rightarrow C_6H_5OH^+$	
66	27.8	13.1 ±0.2	$\rightarrow C_5H_6^+ + CO$	
65	19.3	14.5 ±0.2	$\rightarrow C_5H_5^+ + CO + H$	
63	6.3	19.4 ±0.3	$\rightarrow C_5H_3^+ + ?$	
55	6.3	13.9 ±0.2	$\rightarrow C_3H_3O^+ + ?$	
50	5.3	18.9 ±0.3	$\rightarrow ?$	
40	10.2	14.9 ±0.2	$\rightarrow C_3H_4^+ + ?$	
39	18.0	17.6 ±0.3	$\rightarrow C_3H_3^+ + ?$	
38	12.7	20.4 ±0.9	$\rightarrow C_3H_2^+ + ?$	
29	7.0		$\rightarrow ?$	
27.5	7.7		$\rightarrow C_5H_5^{++} + CO + H$	

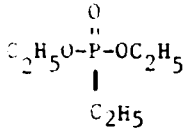
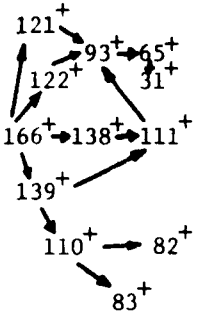
a) Estimated $\Delta H_f[(C_6H_5O)_2POH]$ = -129 kcal/mole.

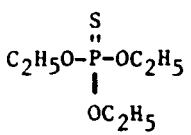
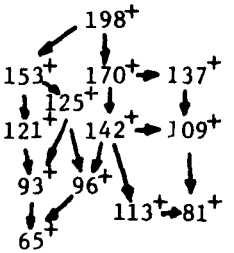
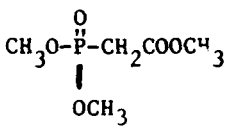
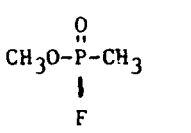
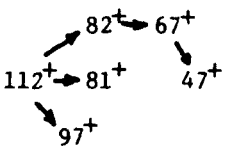
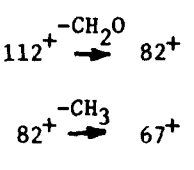
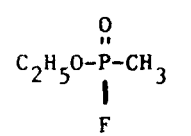
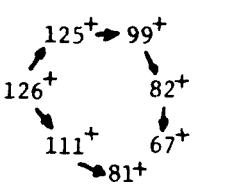
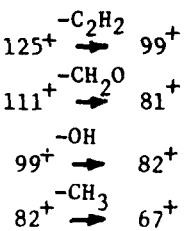
Table 5. Summary of Mass Spectrometric Studies and Findings for Organophosphorus Compounds.

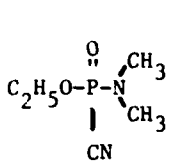
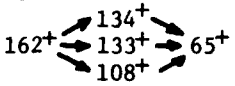
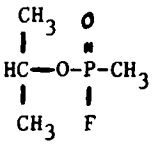
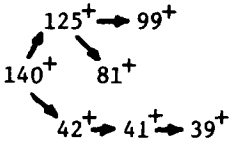
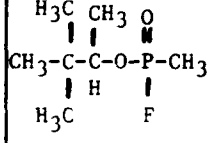
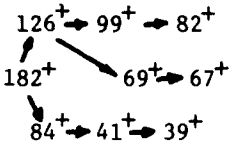
Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
$CH_3-P(Cl)-Cl$	Parent ion is seen; other ions as well; nothing particularly unusual about mass spectrum.	I.P. = 9.49 eV; A.P.'s increase in a rather normal fashion.	$116^+ \rightarrow 101^+ \rightarrow 66^+$; $116^+ \rightarrow 81^+ \rightarrow 45^+$; $116^+ \rightarrow 46^+ \rightarrow 44^+$; $46^+ \rightarrow 15^+$	None observed. (TOF)	Cl^- is the only ion observed.
$Cl-P(=O)-Cl$ $\vert$ Cl	Parent ion is observed. Others about as anticipated.	I.P. = 11.4 eV (EC); A.P.'s increase in a rather normal manner.	$152^+ \rightarrow 117^+ \rightarrow 82^+ \rightarrow 66^+$; $117^+ \rightarrow 47^+$	None observed. (MS-9)	$POCl_2^-$, PO^- and Cl^- were found.
$CH_3-P(=O)-Cl$ $\vert$ Cl	Trimer species was observed; apparently formed as a result of hydrolysis; in "hot source". MW corresponds to: 	I.P. = ca. 11.5 ev (EC); A.P.'s increase in a rather normal manner.	Very difficult to interpret. However, $234^+ \rightarrow 219^+$	Have not been studied.	Have not been studied.

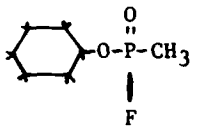
Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
	Parent ion is observed; Others about as expected.	I.P. = 9.95 eV; A.P.'s increase in about the normal manner.		None observed. (TOF)	Cl^- is most intense; also observed $PSCl_2^-$ and PC^- ions.
	Parent ion is observed.	I.P. = 10.48 eV; (10.43 on MS-9); low A.P. for 94^+ (formed by loss of CH_2O). Otherwise, about normal.		$124^+ \xrightarrow{-CH_2O} 94^+$ (RMU-6E)(MS-9) $94^+ \xrightarrow{-CH_3} 79^+$ (RMU-6E)(MS-9) $109^+ \xrightarrow{-CH_2O} 79^+$ (RMU-6E)(MS-9) $93^+ \xrightarrow{-CH_2O} 63^+$ (RMU-6E)(MS-9) $79^+ \xrightarrow{-CH_3OH} 47^+$ (TOF)	Have not been studied.
	Parent ion is observed; remainder of spectrum is about as anticipated.	I.P. = 10.77 eV (10.73 with MS-9); A.P.'s increase in a rather normal manner.		$140^+ \xrightarrow{-CH_2O} 110^+$ (MS-9) $110^+ \xrightarrow{-CH_2O} 80^+$ (MS-9) $109^+ \xrightarrow{-CH_2O} 79^+$ (MS-9) $109^+ \xrightarrow{-CH_4} 93^+$ (TOF)	Have not been studied.

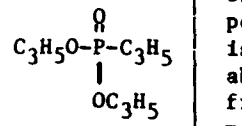
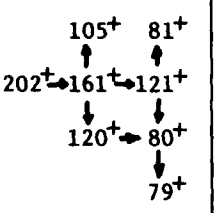
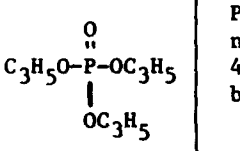
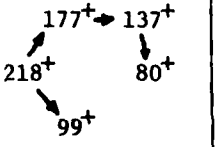
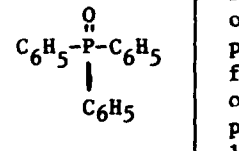
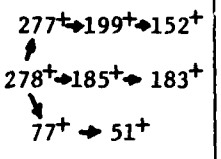
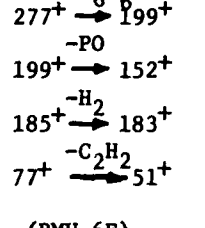
Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
	Parent ion is observed; 93^+ is base peak; some indication of isomerization of ionic species to thiolate structures.	I.P. = 9.2 eV (EC) by TOF; A.P.'s are found to increase in a rather normal manner.	Rather difficult to interpret; however, 	$156^+ \xrightarrow{-CH_2O} 126^+$ $126^+ \xrightarrow{-SH} 93^+$ $125^+ \xrightarrow{-CH_2S} 79^+$ $110^+ \xrightarrow{-CH_2O} 80^+$ $109^+ \xrightarrow{-CH_2O} 79^+$ $93^+ \xrightarrow{-CH_2O} 63^+$ (RMU-6E)	Have not been studied.
	Parent ion is observed.	I.P. = 10.06 eV (10.02 on MS-9); A.P.'s increase in a normal manner.		$182^+ \xrightarrow{-C_2H_3} 155^+$ $181^+ \xrightarrow{-C_2H_4} 153^+$ $167^+ \xrightarrow{-C_2H_4} 139^+$ $182^+ \xrightarrow{-C_2H_4O} 138^+$ $155^+ \xrightarrow{-C_2H_4} 127^+$ $153^+ \xrightarrow{-C_2H_4} 125^+$ $139^+ \xrightarrow{-C_2H_4} 111^+$ $127^+ \xrightarrow{-C_2H_4} 99^+$ $99^+ \xrightarrow{-H_2O} 81^+$ $110^+ \xrightarrow{-C_2H_4} 82^+$ $126^+ \xrightarrow{-C_2H_3O} 83^+$ (MS-9)	Have not been studied.

Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
	Parent ion is observed.	I.P. = 10.39; A.P. of 94^+ is low; C_2H_4 and C_2H_4O are products here.		$166^+ \xrightarrow{-C_2H_3} 139^+$ $166^+ \xrightarrow{-C_2H_4} 138^+$ $166^+ \xrightarrow{-C_2H_4O} 122^+$ $138^+ \xrightarrow{-C_2H_3} 111^+$ $139^+ \xrightarrow{-C_2H_4} 111^+$ $166^+ \xrightarrow{-C_2H_5O} 121^+$ $111^+ \xrightarrow{-H_2O} 93^+$ $109^+ \xrightarrow{-H_2O} 91^+$ $122^+ \xrightarrow{-C_2H_4} 94^+$ $121^+ \xrightarrow{-C_2H_4} 93^+$ $122^+ \xrightarrow{-C_2H_5} 93^+$ $110^+ \xrightarrow{-C_2H_4} 82^+$ $93^+ \xrightarrow{-C_2H_4} 65^+$ $29^+ \xrightarrow{-H_2} 27^+$ $65^+ \xrightarrow{-2(OH)} 31^+$ (RMU-6E)	Have not been studied.

Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
	Parent ion is observed.	I.P. = 8.41 eV; low appearance potential for 93^+; have to postulate S^- formation (?); Otherwise the spectrum is normal.		Studies not completed.	No negative ions are observed.
	No parent ion is observed; 99^+ also is not present; the $P\text{-}CH_2COOCH_3$ bond is rather weak.	I.P. (estimated) = ca. 10.5-11.0 eV; at least four low energy A.P.'s are noted: 110^+, 109^+, 94^+, and 80^+.	Very difficult to interpret.	Studies not completed.	Have not been studied.
EA 1232 	Very small m/q = 99 peak; parent ion is present; $CH_3\text{-}O$ bond is stronger than for alkoxy groups in other G agents.	I.P. = 11.7 eV; A.P. (82^+) = 12.0 ev; A.P.'s of organic fragments not very high, indicating low fragmentation.			Low intensity F^- observed.
EA 1207 	No parent ion; fair-sized $(parent\text{-}H)^+$: also there is a fair-sized $(parent\text{-}CH_3)^+$ ion.	I.P. estimated = ca. 11.5-12.0 eV; A.P. $(99)^+$ is lower than the I.P. (125^+); some processes are quite energetic.	Four ions show maxima: 99^+, 82^+, 111^+ and 125^+; The 99^+ acts as a pseudo-parent. 		Low intensity F^- observed.

Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
GA 	Contains parent ion; many P-contg. ions do not contain CN group; most of the possible combinations of the ions are observed; base peak is m/q = 133; considerable impurities present; m/q = 70 remains unexplained; no m/q=99.	I.P. = 9.44 eV; A.P. (70) = 11.65 eV; A.P.'s of m/q = 44 and 45 are low; they probably come from m/q = 70.	Clastogram including m/q = 70, 44 and 45 does not make much sense; In clastogram without these ions, the parent ion (162^+) behaves as expected; m/q = 133, 108 and 123 show maxima; process suggested is: 	Have not been studied.	Low intensity CN^- observed.
GB 	No parent ion; fair sized $(parent\text{-}CH_3)^+$ ion; very small 67^+ and 82^+; not as much hydrogen rearrangement as with other compounds.	I.P. (estimated) = ca. 12. eV; A.P. of 99^+ is lower than that for 125^+.		Have not been studied.	In Melpar report many negative ions were listed; we have not been able to repeat these data; only observe a low intensity F^- ion.
GD (EA 1210) 	No parent ion; the base peak is 41^+; a large 126^+ ion is $(parent\text{-}C_4H_8)^+$; also a large abundance of organic fragments.	I.P. (estimated) = ca. 12. eV; organic fragments involve highly energetic processes.		Have not been studied.	F^- ion is of low intensity.

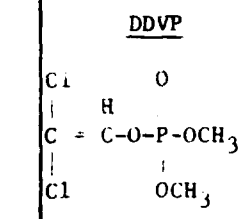
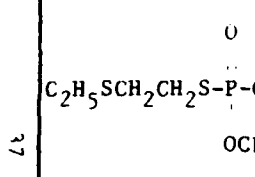
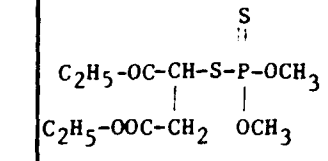
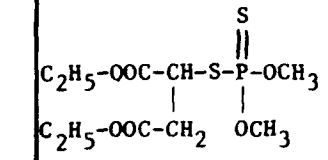
Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
GF 	No parent ion; very weak (R-O) bond; essentially no fragment ions above m/q = 99; no ion observed with cyclohexyl group attached to a P-containing fragment.	I.P. (estimated) = ca. 11. eV; A.P. for 67^+ is low; the A.P.'s of the organic fragments are not very great.	$99^+ \rightarrow 82^+ \rightarrow 67^+$ $180^+ \quad 54^+$ $83^+ \rightarrow 41^+ \rightarrow 39^+$	$99^+ \xrightarrow{-OH} 82^+$ $82^+ \xrightarrow{-CH_3} 67^+$ (TOF)	F^- ion of low intensity.
$C_3H_5O\text{-}P\text{-}OH$ OC_3H_5	Parent ion is not observed; base peak is 41^+, 57^+ is also very abundant; gross fragmentation may be thermal in origin.	A.P. for 121^+ = 10.6 eV; suggests I.P. = 10.0 eV, approximately.	$121^+ \rightarrow 80^+$ $162^+ \quad 81^+ \rightarrow 79^+$ $122^+ \rightarrow 83^+$	Have not been studied.	Have not been studied.
$C_3H_5O\text{-}P\text{-}OC_3H_5$ OC_3H_5	Parent ion is not observed; base peak is 41^+; gross fragmentation may be thermal in origin; sample tends to polymerize in inlet system.	A.P. for 161^+ = 9.15 eV (EC); this suggests I.P. = approximately 8.8 ev; A.P.'s of 79^+, 80^+ and 81^+ are surprisingly low.	105^+ 81^+ $161^+ \rightarrow 121^+$ 80^+ $202^+ \quad 120^+$ 79^+ $41^+ \rightarrow 39^+$	$161^+ \xrightarrow{-C_3H_4} 121^+$ (TOF) Probably many more may be observed with a magnetic instrument.	Have not been studied.

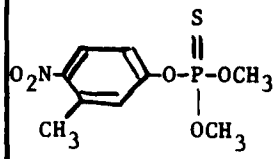
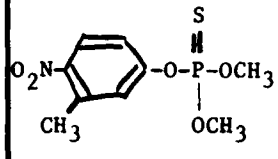
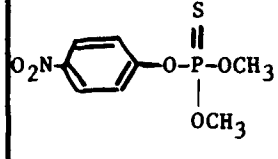
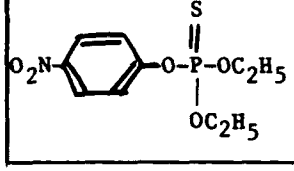
Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
	Parent ion is not observed; base peak is 41^+; 57^+ is also very abundant; gross fragmentation may be thermal in origin.	A.P. for 161^+ = 10.4 ev (EC) this suggests I.P. = approximately 10.0 ev; A.P.'s of 79^+, 80^+ and 81^+ are surprisingly low.		Have not been studied.	Have not been studied.
	Parent ion is not observed; 41^+ is the base peak.	A.P. of 177^+ 10.2 eV (EC); this suggests I.P. = approximately 9.8 eV.		Have not been studied.	Have not been studied.
	Parent ion is observed; base peak is $(P-1)^+$; fragmentation occurs through phenyl and H losses, with and without oxygen loss; P^{++} also seen.	Have not been studied.		 (RMU-6E)	Have not been studied.

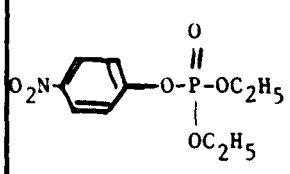
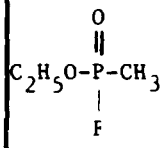
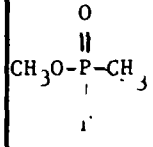
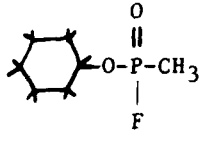
Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
$C_6H_5O-P(OH)-OC_6H_5$	Parent ion is not observed; base peak is 94^+ ion; mass spectrum is very suggestive of thermal decomposition to yield mostly phenol.	A.P. of 94^+ = 8.5 eV (EC): I.P. of phenol = 8.50 eV by P.I.; thus, 94^+ is $C_6H_5OH^+$, apparently formed by simple ionization of phenol.	Not too meaningful, since apparently largely due to phenol from thermal decomposition.	Have not been studied.	Have not been studied.
$C_6H_5O-P(OC_6H_5)-OC_6H_5$	Parent ion is not observed; base peak is 94^+ ion; mass spectrum is very suggestive of thermal decomposition to yield mostly phenol; small amounts of 149^+, 156^+, and 167^+ observed.	A.P. of 94^+ = 8.8 ± 0.2 eV (EC): I.P. of phenol = 8.50 eV by P.I.; thus, 94^+ is $C_6H_5OH^+$, apparently formed by simple ionization of phenol.	Not too meaningful, since apparently largely due to phenol from thermal decomposition.	Have not been studied.	 most intense at m/q = 93; also $(C_6H_5O)PO^-$ at m/q = 233 observed; others at m/q = 133 and 217.

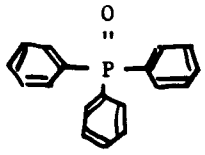
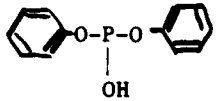
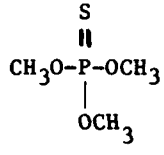
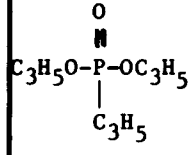
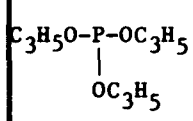
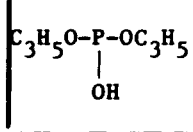
Formula	Mass Spectra	Energetics	Clastograms	Metastables	Negative Ions
$C_6H_5NO_2$	Both parent positive and negative ions observed; base peak in negative ion spectrum is NO_2^-; base peak in positive ion spectrum is $C_6H_5^+$.	Have not been studied.	Have not been studied.	$123^+ \xrightarrow{-NO} 93^+$ $123^+ \xrightarrow{-NO_2} 77^+$ $123^+ \xrightarrow{-HNO_2} 76^+$ $107^+ \xrightarrow{-N} 93^+$ $107^+ \xrightarrow{-NO} 77^+$ $93^+ \xrightarrow{-CO} 65^+$ $65^+ \xrightarrow{-C_2H_2} 39^+$ $39^+ \xrightarrow{-H_2} 37^+$ $77^+ \xrightarrow{-C_2H_2} 51^+$ $75^+ \xrightarrow{-C_2} 51^+$ $123^+ \xrightarrow{-C_6H_5O} 30^+$ $77^+ \xrightarrow{-C_4H_5} 24^+$ $77^+ \xrightarrow{-C_4H_2} 27^+$ $77^+ \xrightarrow{-C_2H_3} 50^+$ $76^+ \xrightarrow{-C_2H_2} 50^+$ (RMU-6E)	NO_2^- is very intense; $C_6H_5O^-$ also observed along with $C_6H_5NO_2^-$ in rather large yields.

Table 9 A Short Catalog of Mass Spectra of Organophosphorus and Related Compounds.

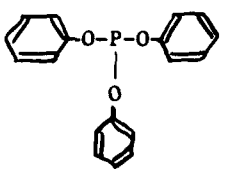
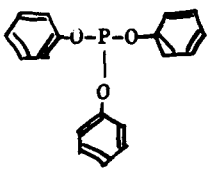
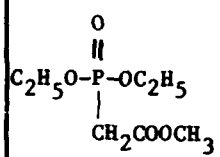
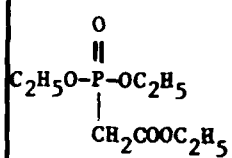
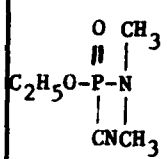
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
DDVP 	109 102.	185,7 33.	79 27.	145,7 11	220,2,4 11.	47 11	60 2				220+ 222+ 224+	10
Metasystox 	88 112	60 59.	109,10,11 20.	142,3,4 19.	79 11.	59 9.	75 7.	230,1,2 7.	45 6.	125 5.	230+ 231+ 232+	10
Malathion 	173 111.	127,8 91.	125,6 89.	93 73.	158,9,60 63.	99 30.	55 28.	143 20.	43 19.	79 18.	330+ 331+ 332	10
Malathion 	29 49.	173 120.	127,8 97.	125,6 61.	93 54.	158,9,60 52.	99 22.	27 17.	63 16.	47 16.	330+ 331+ 332	11

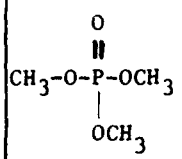
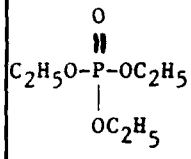
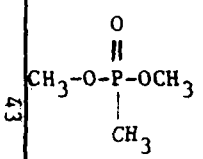
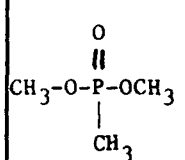
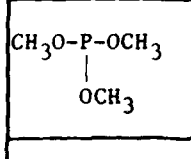
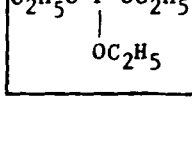
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
Folithion 	109,10,11 / 121.	277,8,9 / 106.	125,6 / 102.	260,1,2 / 51.	79 / 38.	47 / 36.	93 / 31.	63 / 19.	89 / 10.	121 / 7.	277+ 278+ 279	10
Folithion 	125,6,7 / 121.	109,10,11 / 111.	277,8 / 95.	260,1,2 / 60.	79 / 47.	47 / 41.	93 / 32.	63 / 34.	51 / 20.	52 / 17.	277+ 278	11
Methyl Parathion 	109,10 / 106.	125 / 70.	263 / 52.	79 / 29.	44 / 29.	47 / 26.	63 / 26.	93 / 25.	15 / 15.	43 / 14.	263	11
Parathion 	109 / 100.	29 / 85.	291 / 84.	97 / 72.	139 / 62.	137 / 60.	40 / 49.	125 / 45.	155 / 34.	43 / 30.	291	11

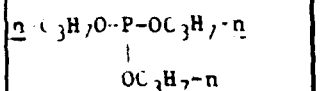
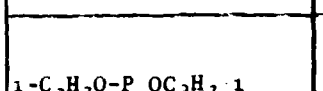
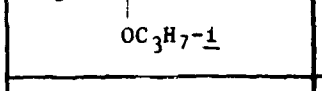
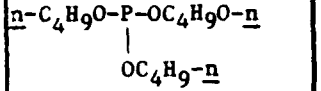
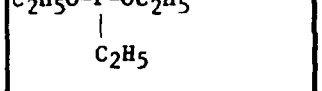
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M W	Ref.
Paraoxon 	109,110 / 114	29 / 56	149 / 47	81 / 46	275 / 40	139 / 37	65 / 32	75 / 25	63 / 25	64 / 25	275	11
EA 1207 	99 / 100	81 / 31.6	82 / 17.2	111 / 15.4	29 / 10.1	47 / 6.8	83 / 6.2	67 / 6.0	27 / 5.6	15 / 4.6	126	12
EA 1232 	82 / 100.0	67 / 43.8	81 / 24.6	47 / 13.9	15 / 13.4	83 / 12.3	112 / 12.2	97 / 10.9	29 / 7.3	31 / 5.9	112	12
GF 	99 / 100.0	67 / 28.7	41 / 26.8	54 / 24.7	39 / 20.3	27 / 17.2	29 / 16.4	81 / 13.8	55 / 12.8	82 / 11.7	180	12

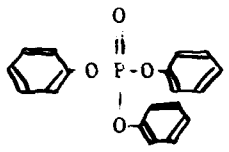
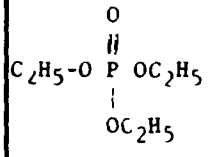
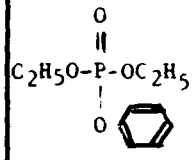
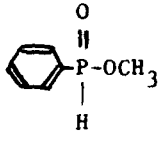
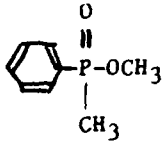
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
	277,8 120.0	278,9 31.7	201,2 24,0	199,200 22,0	183,4 17,7	77 16,6	185,6 12,9	152,3 10.6	51 8.4	154,5 5.3	278	12
	93,4 119,5	66 28,2	65 19,7	39 18,3	40 10,3	27 7,8	38 7,7	29 7,0	63 6,5	55 6,3	234	12
	93,4,5 117.	156,8 31,	83,5 21,5	126,8 21,3	47 19,6	56 17,5	63 17,0	31, 14,7	29 11,0	79 11,0	156+ 158	12
	41 100.0	57 62.1	39 49.2	79 37,2	81 27.5	80 26,5	58 16.5	121 13,3	55 12.5	67 11,3	202	12
	41 100.0	39 21.3	81 14.8	79 13.1	80 9.5	57 9,0	42 6,3	55 5.1	54 4.5	161 4.3	202	12
	41 100.0	57 69.8	39 54,5	121 36.5	58 25.0	80 22.8	42 13.5	29 14.7	27. 12.9	55 10.9	162	12

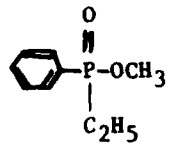
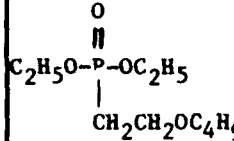
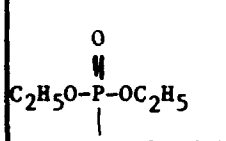
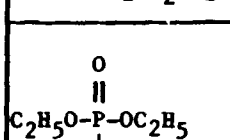
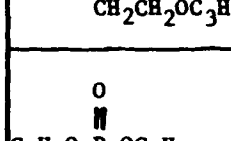
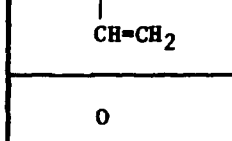
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
$CH_3-P(Cl)_2$	101 100.0	43 84.2	116 70.2	44 22.4	81 21.6	46 9.0	35 8.1	31 5.5	15 5.3	66 5.3	116+ 118+ 120	12
$Cl-P(=O)(Cl)_2$	117 100.0	154 44.2	47 37.3	35 23.7	31 12.6	66 6.9	82 4.1	101 2.0			154+ 156+ 158	12
$CH_3-P(=S)(Cl)_2$	113 100.0	148 61.5	45 47.5	77 34.5	63 32.5	44 16.2	31 15.6	15 13.7	133 10.5	43 8.0	148+ 150	12
$C_2H_5O-P(=O)(OC_2H_5)_2$	111 100.0	93 84.1	31 70.1	65 68.6	139 53.3	81 36.9	166 34.4	45 32.9	82 28.8	78 27.9	166	12
$C_2H_5O-P(=S)(OC_2H_5)_2$	29 100.0	65 64.0	121 59.3	27 56.4	97 56.4	198 47.7	93 45.3	45 37.2	115 32.6	109 30.2	198+ 200	12
$C_2H_5O-P(=O)(CH(CH_3)COOC_2H_5)(OC_2H_5)$	109 100.	193 94.	165 82.	137 73.	29 66.	81 65.	155 50.	99 45.	138 40.	27 36.	238	13

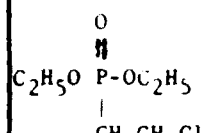
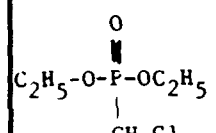
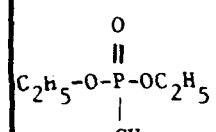
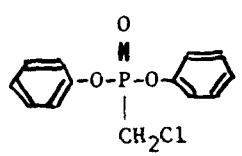
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
	94 / 100.0	66 / 32.0	65 / 23.7	39 / 23.3	95	93	43	54	38	62	310	12
	94 / 100.0	217 / 55.	77 / 46.	43 / 35.	66 / 22.	65 / 21.	39 / 18.	51 / 15.	310 / 14.	153 / 14.	310	14
	183 / 100.	123 / 73.	74 / 48.	155 / 46.	137 / 40.	95 / 34.	109 / 33.	29 / 30.	81 / 29.	42 / 28.	210	13
	123 / 100.	197 / 72.	179 / 71.	151 / 63.	29 / 63.	109 / 52.	152 / 48.	81 / 47.	42 / 41.	88 / 41.	224	13
GA 	43 / 100.0	44 / 76.3	70 / 52.6	42 / 45.8	15 / 17.0	133 / 14.1	45 / 12.4	47 / 10.2	162 / 9.0	18 / 6.8	162	12

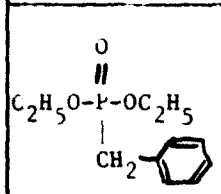
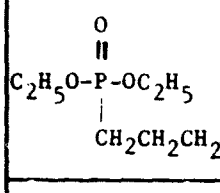
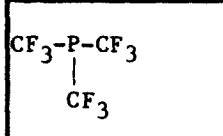
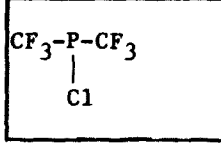
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
	15/100.0	110/86.3	29/36 6	109/28.0	79/26.6	95/22.1	47/22.0	80/17.3	31/16.7	140/14.0	140	15
	99/100.0	81/78.3	155/68.0	109/49.3	127/47.0	82/47.0	45/45.4	182/26.0	29/25.0	125/19.7	182	15
	79/100.0	94/88.0	15/82.5	47/47.6	29/39.8	93/38.6	109/28.3	80/21.7	31/16.9	124/15.7	124	15
	79/100.0	94/81.5	47/37.6	15/34.7	109/33.0	93/31.6	63/20.8	124/13.6	29/12.4	31/10.7	124	16
	93/100.0	63/64.	109/57.	15/42.	79/41.	94/36.	47/36.	124/36.	80/23.	31/20.	124	14
	65/100.0	82/94.	83/76.	111/76.	29/56.	139/52.	93/46.	81/43.	121/37.	166/36.	166	14

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref
	83 / 100.	43 / 42.	82 / 34.	65 / 27.	125 / 21.	41 / 17.	107 / 14.	27 / 12.	167 / 11.	149 / 10.	208	14
	82 / 100.	124 / 62.	43 / 62.	109 / 49.	83 / 48.	107 / 47.	65 / 30.	41 / 29.	45 / 20.	27 / 15.	208	14
	83 / 100.	57 / 30.	138 / 25.	41 / 20.	29 / 16.5	56 / 13.	195 / 9.	31 / 6.	177 / 5.	65 / 5.	250	14
	111 / 100.0	93 / 69.	139 / 47.	65 / 44.	29 / 35.	109 / 34.	138 / 30.	82 / 28.	81 / 23.	83 / 21.	166	14
	83 / 100.	109 / 85.	45 / 39.	43 / 35.	41 / 19.	27 / 12.	59 / 11.	39 / 10.	65 / 10.	42 / 8.	166	14
	97 / 100.	123 / 65.	45 / 31.	43 / 27.	79 / 24.	41 / 22.	28 / 18.	39 / 15.	27 / 10.	80 / 9.	180	14

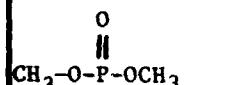
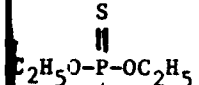
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W	Ref
	326 100.	325 81	94 74.	77 77	170 40.	169 26	233 25	215 22	168 14.	232 13	326	18
	155 100.	99 93.	127 56.	81 43.	109 41.	82 31.	29 19	125 18	111 15.	45 15.	182	17
	94 100.	104. 78.	65 27.	77 24.	230 23.	81 21.	174 17.	187 14.	66 13.	202 12.	230	18
	77 100.	79 96.	51 88.	47 74.	78 69.	155 56.	156 42.	50 35.	126 33.	141 22.	156	19
	155 100.	77 65	47 42.	170 42.	51 28.	169 27.	140 26.	91 24.	125 20.	139 8.	170	19

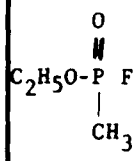
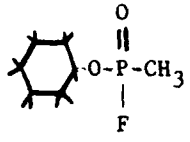
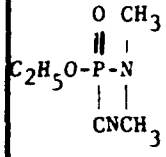
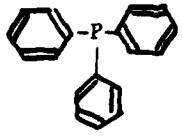
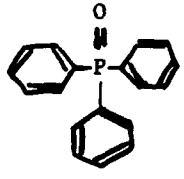
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
	155/100.	77/70.	51/40.	47/33.	184/27.	78/14.	50/13.	79/12.	91/12.	141/9.	184	19
	65/100.	109/96.	181/88.	138/77.	111/65.	166/58.	81/54.	137/50.	57/35.	139/34.	238	20, 21
	125/100.	81/79.	111/75.	82/60.	138/55.	29/44.	65/38.	109/35.	153/34.	137/29.	210	21
	125/100.	81/88.	111/80.	138/71.	82/55.	65/47.	137/46.	109/41.	43/39.	139/33.	224	21
	91/100.	109-/87.	109+/86.	27/61.	29/51.	137/47.	136/36.	119/19.	45/18.	110/8.	164	22
	109/100.	79/28.	15/13.	158+60/8.	49+51/5.	31/4.	127+9/2.				158+160	22

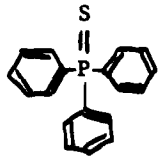
Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W	Ref.
	138 100.	145+7 80.	109 62	127+9 60.	29 49.	173+5 44	91 37.	45 25.	137 16.	110 15.	200+ 202	22
	109 100.	159+161 66.	29 43.	131+3 41.	91 40.	137 35.	45 26.	113+5 24.	93 15.	141+3 11.	186+ 188	22
	97 100.	125 90.	79 77.	124 26.	107 26.	45 10.	109 9.	93 8.	137 7.	29 6.	152	22
	125+7 100.	77 83.	282+4 66.	170 30.	165 30.	94 24.	215 17.	93 5.	49+51 3.		282+ 284	22
	91 100.	167 53.	324 51.	77 11.	94 6.	93 2.					324	22

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
	91 100.	109 25.	228 17.	172 14.	29 14.	200 6.	93 5.	45 3.	173 1.	137 1.	228	22
	125 100.	152 87.	111 78.	139 70.	138 50.	65 46.	97 44.	81 38.	82 26.	109 34.	194	21
CF_3PCl_2	101+3+5 100.0	170+2+4 25.4	85+7 20.2	66+8 10.5	69^+ 7.9	69^- 6.3	151+3+5 3.1	50 1.8	31^+ 1.6	31^- 1.4	170+ 172+ 174	23
CF_3PF_2	69^+ 100.0	69^- 58.4	31 11.2	138 9.0	119 6.9	50^+ 4.9	50^- 3.9				138	23
CF_3PH_2	102 100.	69^+ 100.	69^- 49.5	51^+ 16.8	32 11.7	52 12.8	83 11.4	50 11.0	51^- 8.4	33 8.1	102	23
	69^+ 100.	69^- 24.8	131 16.4	219 9.0	100 8.8	31 7.8	150 5.7	238 4.8	50 4.6	81 3.6	238	23
	69^+ 100.	85+7 83.0	204+6 42.6	69^- 31.8	116+8 27.8	66+8 9.2	31 7.9	50 5.3	135+7 4.6	100 3.9	204+ 206	23

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
CF_3 P-CF_3 | F	69 / 100.	69^+ / 88.5	31 / 27.6	100 / 24.6	188 / 24.2	119 / 15.5	50^- / 10.6	81 / 4.5	169 / 3.9	50^+ / 3.9	188	23
CF_3-P-CF_3 | H	69^+ / 100.	69^- / 31.8	170 / 25.6	82 / 24.4	151 / 11.2	101 / 10.4	51^+ / 10.4	51^- / 7.9	31 / 7.9	50 / 4.8	170	23
F-P-F | F	69 / 100.	88 / 43.4	50 / 9.5	31 / 5.7	19 / 4.6						88	24
O ‖ F-P-F | F	85 / 100.0	104 / 87.0	69 / 17.7	50 / 4.4	31 / 3.6	47 / 2.2	88 / 1.	66 / 1.	19 / 1.		104	24
O ‖ Cl-P-Cl | Cl	117+9+21 / 100.0	152+4+6 / 44.2	47 / 37.3	35+7 / 23.7	31 / 12.6	66+8 / 6.9	82+4 / 4.1				152+ 154+ 156	24
O ‖ Br-P-Br | Br	205+7+9 / 100.0	47 / 66.6	284+6+ 288+90 / 43.4	79+81 / 37.	126+8 / 15.1	31 / 9.6	110+12 / 6.2		189+91+ 193 / 0.9		284+ 286+ 288+ 290	24

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
	120+1+2 100.	32+3+4 73.5	101+2+3 27.1	69 26.0	88 6.9	50 6.8	63+4+5 2.4				120+ 121+ 122	24
	101+3+5 100.	45 84.2	116+8+20 70.2	44 22.4	81+83 21.6	35+7 8.1	31 5.5	66+8 5.3	15 5.3	80+2 4.6	116+ 118+ 120	24
	113+4+5 100.	148+50+ 152 61.5	45 47.5	77+8+9 34.5	63+4+5 32.5	44 16.2	31 15.6	15 13.7	133+5+7 10.5	43 8.0	148+ 150+ 152	24
	79 100.	80 91.8	15 49.4	47 45.9	31 36.7	29 33.6	110 19.3	94 16.3	109 15.3	95 14.3	182	24
	29 100.	27 59.2	121 56.3	65 54.3	97 51.8	198 42.7	93 41.4	45 33.5	109 26.4	81 23.3	198	24
	82 100.0	67 43.8	81 24.6	47 13.9	15 13.4	83 12.3	112 12.2	97 10.9	29 7.3	31 5.9	112	24

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref
	99 / 100.0	81 / 31.6	82 / 17.2	111 / 15.4	29 / 10.1	47 / 6.8	83 / 6.2	67 / 6.0	27 / 5.6	15 / 4.6	126	24
	99 / 100.0	67 / 28.3	41 / 26.2	54 / 24.0	39 / 20.0	27 / 17.2	29 / 16.4	81 / 12.7	55 / 11.5	82 / 11.3	180	24
	133 / 100.0	47 / 73.4	162 / 66.4	108 / 41.8	117 / 34.9	107 / 32.1	134 / 31.0	106 / 31.0	65 / 29.0	135 / 20.3	162	24
	262 / 100.	183 / 44.	263 / 23.	108 / 21.	184 / 10.	107 / 10.	185 / 7.	152 / 5.	51 / 5.	154 / 3.5	262	25
	277 / 100.	278 / 53.	77 / 30.	201 / 19.	183 / 14.	199 / 13.	51 / 10.	185 / 9.	47 / 8.	152 / 7.	278	25

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
	294 100.	183 54.	185 47.	293 40.	262 32.	295 18.	43 15.	152 13.	263 11.	139 9.	294	25
H-P-H | H	32 100.	34 75.2	31 24.1	33 21.8							34	26
H-P-H | H	32 100.0	34 85.0	33 27.6	31 26.4							34	27
H-P-H | H	32 100.0	34 68.0	31 40.2	33 25.4							34	28
H-P-H | H	32 100.0	34 89.4	33 26.2	31 18.7							34	29
H-P-H | H	34 100.0	33 33.1	31 32.1	32 12.7							34	30
H-P-P-H | | H H	32 100.0	66 74.6	31 65.3	62 57.3	34 52.0	64 50.7	63 44.0	33 21.3	65 9.3		66	26

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
H-P-P-H | | H H	62 100.0	66 8.4	64 5.8	63 5.6	33 5.1	65 1.3	31 1.2	34 0.4	32 0.2		66	28
F | F-P-F /\\ F F	107 100.	88 21.0	50 10.3	69 6.0	31 3.2	19 2.0					126	26
Cl-P-Cl | Cl	101+3+5 100.	35+7 48.3	136+8+ 140 31.5	66+8 29.2	31 20.9						136+ 138+ 140	26
O ‖ Cl-P-Cl | Cl	117+9+ 121 100.0	47 87.5	35+7 49.5	152+4+6 37.8	31 21.0	66+8 20.8	82+4 7.4	101+3+5 6.2			152+ 154+ 156	26
S ‖ Cl-P-Cl | Cl	133+4+ 135+6+7 100.0	168+9+ 170+1+2+4 51.3	63 47.1	31 27.3	35+7 25.5	101+3+5 18.5	66+8 11.2	98+9+100 8.0			168+9+ 170+1+2 +174	26
CH_3-P-CH_3 | CH_3	61 100.0	76 78.2	59 66.2	45 56.4	57 32.8	75 20.0	15 15.2	44 12.1	58 11.8	47 11.4	76	26

Formula of Compound	1st	2nd	3rd	4th	5th	6th	7th	8th	9th	10th	M.W.	Ref.
$C_2H_5-P(H)(C_2H_5)$	62 / 100.0	90 / 75.6	61 / 73.0	118 / 52.7	57 / 46.9	59 / 45.7	45 / 29.0	103 / 22.8	58 / 20.7	75 / 19.5	118	26
$CH_3-P(H)-H$	46 / 100.0	15 / 79.7	48 / 63.7	45 / 56.6	44 / 42.6	47 / 15.8	43 / 8.3				48	26
$C_2H_5-P(H)-H$	62 / 100.0	34 / 53.2	57 / 50.7	58 / 35.2	45 / 35.2	43 / 25.3	47 / 19.7	59 / 19.4	60 / 18.7	44 / 14.4	62	26
$CH_3-O-P(OCH_3)-OCH_3$	93 / 100.0	15 / 93.6	63 / 51.7	109 / 48.9	47 / 43.0	124 / 38.7	94 / 28.3	79 / 17.0	29 / 16.5	45 / 13.9	124	26
$C_2H_5O-P(OC_2H_5)-OC_2H_5$	29 / 100.0	65 / 68.1	82 / 62.9	27 / 61.3	83 / 35.0	81 / 31.1	111 / 23.1	47 / 20.1	139 / 17.7	79 / 17.1	166	26
$CH_3O-P(OH)-OCH_3$	79 / 100.0	80 / 91.8	47 / 54.6	15 / 51.5	31 / 41.6	29 / 25.8	49 / 18.9	110 / 18.4	95 / 15.9	48 / 12.7	110	26
$C_2H_5C-P(OH)-OC_2H_5$	111 / 100.0	83 / 99.4	65 / 86.5	45 / 79.7	29 / 65.3	31 / 59.3	27 / 49.9	93 / 31.5	66 / 31.3	110 / 19.0	138	26

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13. ABSTRACT

Mass spectrometric studies of trimethyl phosphorothionate, triphenyl phosphite, diallyl phosphite, diallyl allylphosphonate, triallyl phosphite, triallyl phosphate, and diphenyl phosphite were made to obtain more fundamental information for use in determining the ionization and dissociative ionization processes that occur upon electron impact with these organophosphorus compounds. Mass spectra, appearance potentials, clastograms, and metastable transitions obtained in these studies are reported. These results are discussed in relation to a better understanding of the fragmentation pathways in this class of compounds. A summary table of all results determined under this contract is included, and a brief catalog, with literature references, of the mass spectra of organophosphorus compounds summarizes much of the available literature. Future correlative activities are suggested.

14. KEY WORDS:

- Appearance potentials
- Biologically-active compounds
- Clastograms
- Diallyl allyl-phosphonate
- Diallyl phosphite
- Diphenyl phosphite
- Fragmentation processes
- G agents
- Mass Spectra
- Metastable transitions
- Organophosphorus compounds
- Triallyl phosphate
- Triallyl phosphite
- Trimethyl phosphorothionate
- Triphenyl phosphite

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